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PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

***** Welcome to STN International *****

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
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STN Express with Discover!
NEWS 4 OCT 28 KOREAPAT now available on STN
NEWS 5 NOV 30 PHAR reloaded with additional data
NEWS 6 DEC 01 LISA now available on STN
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NEWS 8 DEC 15 MEDLINE update schedule for December 2004
NEWS 9 DEC 17 ELCOM reloaded; updating to resume; current-awareness
alerts (SDIs) affected
NEWS 10 DEC 17 COMPUAB reloaded; updating to resume; current-awareness
alerts (SDIs) affected
NEWS 11 DEC 17 SOLIDSTATE reloaded; updating to resume; current-awareness
alerts (SDIs) affected
NEWS 12 DEC 17 CERAB reloaded; updating to resume; current-awareness
alerts (SDIs) affected
NEWS 13 DEC 17 THREE NEW FIELDS ADDED TO IFIPAT/IFIUDE/IFICDB
NEWS 14 DEC 30 EPFULL: New patent full text database to be available on STN
NEWS 15 DEC 30 CAPLUS - PATENT COVERAGE EXPANDED
NEWS 16 JAN 03 No connect-hour charges in EPFULL during January and
February 2005
NEWS 17 FEB 25 CA/CAPLUS - Russian Agency for Patents and Trademarks
(ROSPATENT) added to list of core patent offices covered
NEWS 18 FEB 10 STN Patent Forums to be held in March 2005
NEWS 19 FEB 16 STN User Update to be held in conjunction with the 229th ACS
National Meeting on March 13, 2005
NEWS 20 FEB 28 PATDPAFULL - New display fields provide for legal status
data from INPADOC
NEWS 21 FEB 28 BABS - Current-awareness alerts (SDIs) available
NEWS 22 FEB 28 MEDLINE/LMEDLINE reloaded

NEWS EXPRESS JANUARY 10 CURRENT WINDOWS VERSION IS V7.01a, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 10 JANUARY 2005

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.

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of commercial gateways or other similar uses is prohibited and may
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***** STN Columbus *****

FILE 'HOME' ENTERED AT 08:17:57 ON 02 MAR 2005

=> FIL STNGUIDE

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

0.21

0.21.

FILE 'STNGUIDE' ENTERED AT 08:18:02 ON 02 MAR 2005

USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT

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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Feb 25, 2005 (20050225/UP).

=> FIL HOME

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

0.06

0.27

FILE 'HOME' ENTERED AT 08:18:07 ON 02 MAR 2005

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

0.21

0.48

FILE 'REGISTRY' ENTERED AT 08:18:15 ON 02 MAR 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 28 FEB 2005 HIGHEST RN 839671-97-5

DICTIONARY FILE UPDATES: 28 FEB 2005 HIGHEST RN 839671-97-5

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

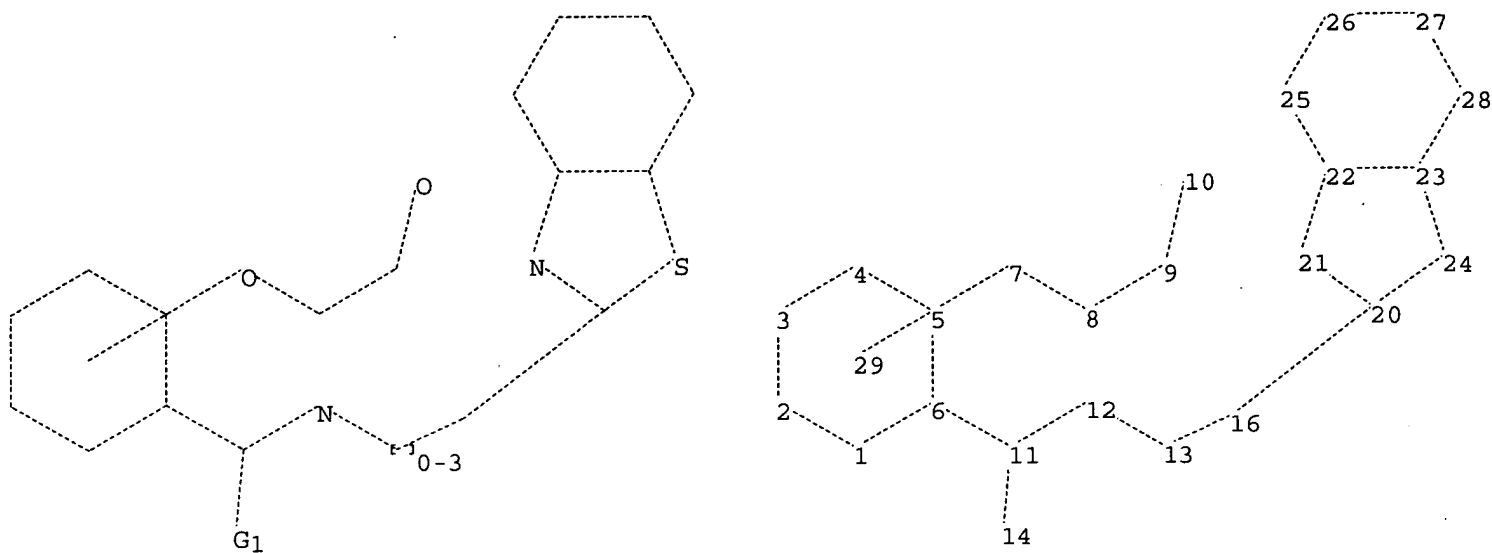
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:

<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10662135.str



chain nodes :

7 8 9 10 11 12 13 14 16

ring nodes :

1 2 3 4 5 6 20 21 22 23 24 25 26 27 28

chain bonds :

6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 16-20

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 16-20 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28

isolated ring systems :

containing 1 : 20 :

G1:O,S,N

Match level :

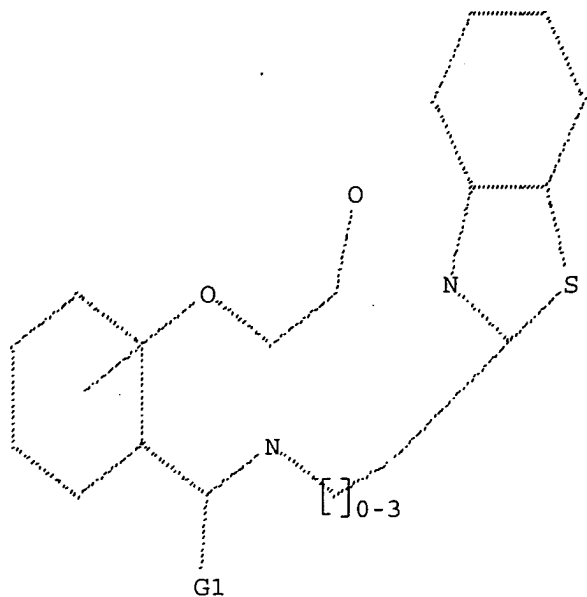
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS
12:CLASS 13:CLASS 14:CLASS 16:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom 29:CLASS

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR

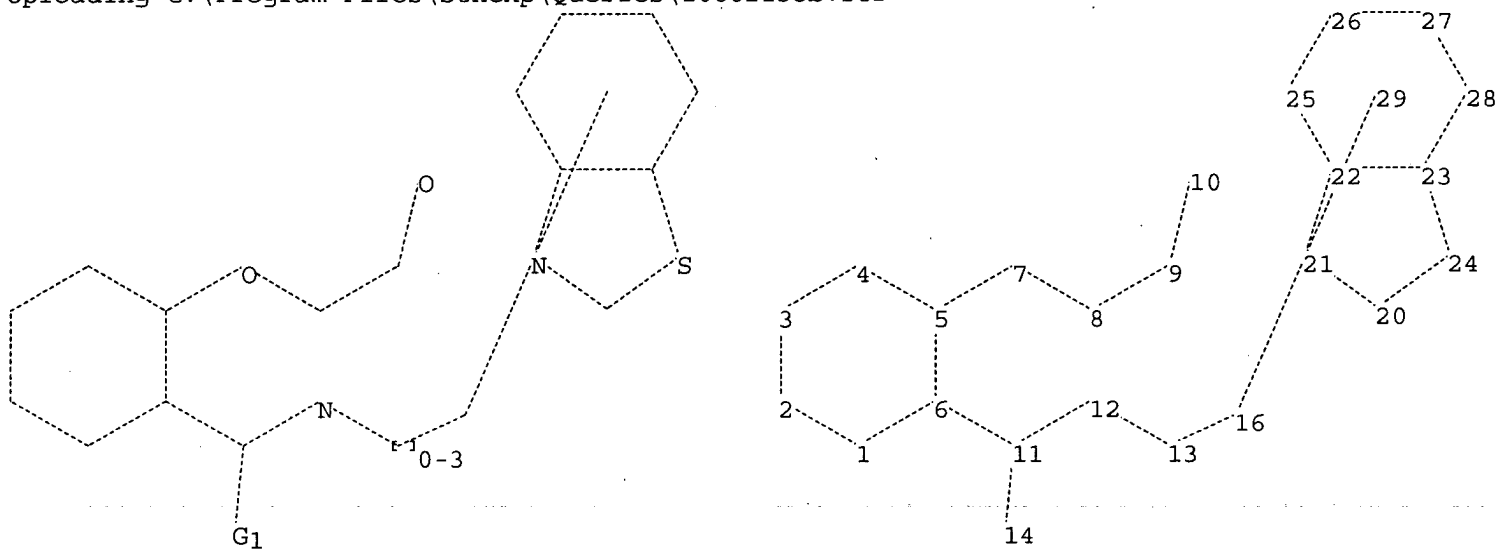


G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

=>

Uploading C:\Program Files\Stnexp\Queries\10662135b.str



chain nodes :

7 8 9 10 11 12 13 14 16

ring nodes :

1 2 3 4 5 6 20 21 22 23 24 25 26 27 28

chain bonds :

5-7 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28

isolated ring systems :

containing 1 : 20 :

G1:O,S,N

Match level :

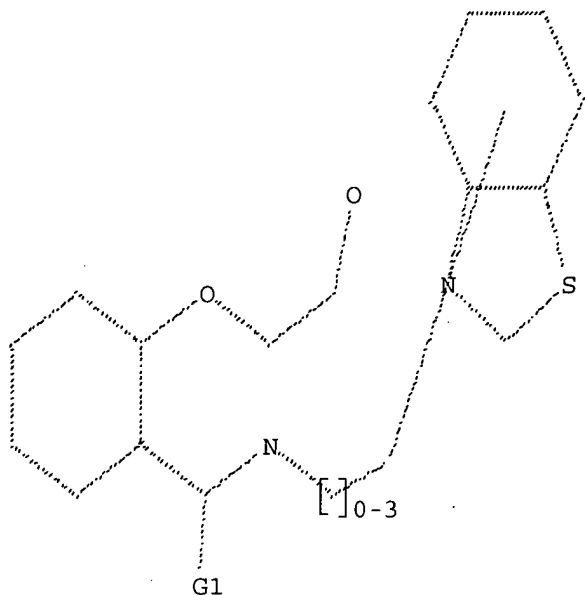
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12:CLASS 13:CLASS 14:CLASS 16:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom 29:CLASS

L2 STRUCTURE UPLOADED

=> d

L2 HAS NO ANSWERS

L2 STR



G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 08:19:45 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA SSS SAM L1

=> s l1 full

FULL SEARCH INITIATED 08:19:50 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS 8 ANSWERS
SEARCH TIME: 00.00.01

L4 8 SEA SSS FUL L1

=> s l2

SAMPLE SEARCH INITIATED 08:19:54 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 2 TO 124
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L2

=> s 12 full
FULL SEARCH INITIATED 08:19:57 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 39 TO ITERATE

100.0% PROCESSED 39 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

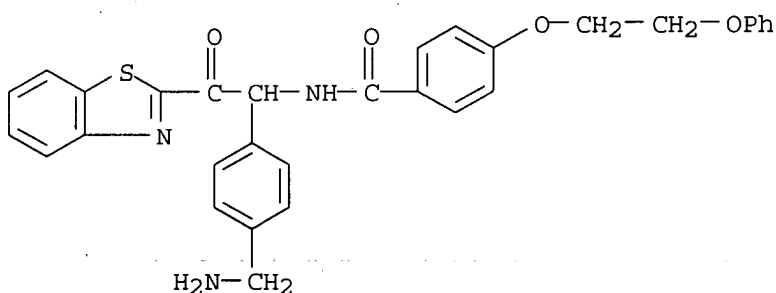
L6 0 SEA SSS FUL L2

=> s 14 and caplus/lc
44163940 CAPLUS/LC
L7 7 L4 AND CAPLUS/LC

=> s 14 not 17
L8 1 L4 NOT L7

=> d

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN 334989-04-7 REGISTRY
CN Benzamide, N-[1-[4-(aminomethyl)phenyl]-2-(2-benzothiazolyl)-2-oxoethyl]-4-(2-phenoxyethoxy) - (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C31 H27 N3 O4 S
CI COM
SR CA



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

| | | |
|----------------------|------------|---------|
| => fil caplus | | |
| COST IN U.S. DOLLARS | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| FULL ESTIMATED COST | 329.96 | 330.44 |

FILE 'CAPLUS' ENTERED AT 08:20:51 ON 02 MAR 2005
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FILE COVERS 1907 - 2 Mar 2005 VOL 142 ISS 10
FILE LAST UPDATED: 1 Mar 2005 (20050301/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 17
L9

3 L7

=> d ibib abs hitstr l9 1-3

L9 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2004:819180 CAPLUS

DOCUMENT NUMBER: 141:374428

TITLE: Design and synthesis of highly potent and selective (2-arylcarbamoyl-phenoxy)-acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications

AUTHOR(S): Van Zandt, Michael C.; Sibley, Evelyn O.; McCann, Erin E.; Combs, Kerry J.; Flam, Brenda; Sawicki, Diane R.; Sabetta, Al; Carrington, Anne; Sredy, Janet; Howard, Eduardo; Mitschler, Andre; Podjarny, Alberto D.

CORPORATE SOURCE: The Institute for Diabetes Discovery, LLC, Branford, CT, 06405, USA

SOURCE: Bioorganic & Medicinal Chemistry (2004), 12(21), 5661-5675

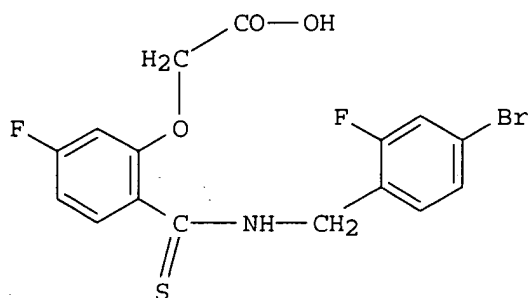
CODEN: BMECEP; ISSN: 0968-0896

PUBLISHER: Elsevier Ltd.

DOCUMENT TYPE: Journal

LANGUAGE: English

GI



AB Recent efforts to identify treatments for chronic diabetic complications have resulted in the discovery of a novel series of highly potent and selective (2-arylcarbamoyl-phenoxy)-acetic acid aldose reductase inhibitors. The compound class features a core template that utilizes an intramol. hydrogen bond to position the key structural elements of the pharmacophore in a conformation, which promotes a high binding affinity. The lead candidate, I, 5-fluoro-2-(4-bromo-2-fluorobenzylthiocarbamoyl)-phenoxyacetic acid, inhibits aldose reductase with an IC₅₀ of 30 nM, while being 1100 times less active against aldehyde reductase, a related enzyme involved in the detoxification of reactive aldehydes. In addition, I lowers nerve sorbitol levels with an ED₅₀ of 31 mg/kg/d po in the 4-day STZ-induced diabetic rat model.

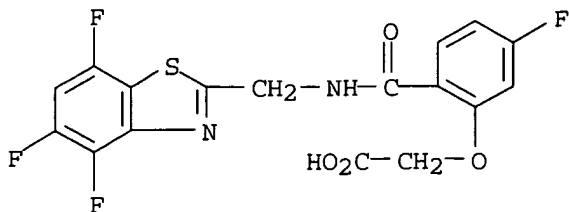
IT 314297-79-5P 314297-80-8P 314297-81-9P
314297-82-0P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(design and synthesis of (arylcarbamoylphenoxy)acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications)

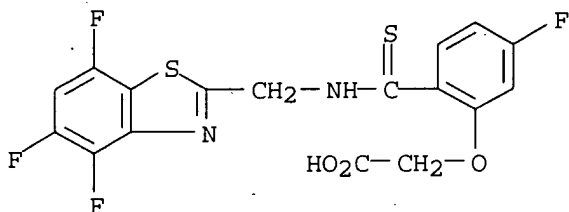
RN 314297-79-5 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



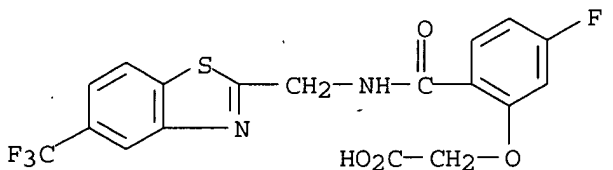
RN 314297-80-8 CAPLUS

CN Acetic acid, [5-fluoro-2-[thioxo[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]- (9CI) (CA INDEX NAME)



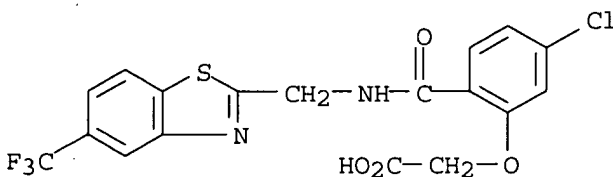
RN 314297-81-9 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



RN 314297-82-0 CAPLUS

CN Acetic acid, [5-chloro-2-[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



IT 314298-38-9P

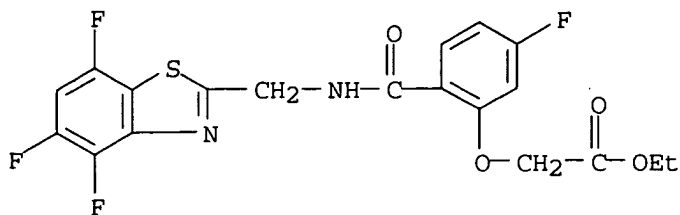
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(design and synthesis of (arylcarbamoylphenoxy)acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications)

RN 314298-38-9 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]-, ethyl ester (9CI) (CA

INDEX NAME)



REFERENCE COUNT: 44 THERE ARE 44 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L9 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2001:283939 CAPLUS

DOCUMENT NUMBER: 134:311433

TITLE: Preparation of (hetero)arylmethylamines as tryptase inhibitors

INVENTOR(S): Lively, Sarah Elizabeth; Waszkowycz, Bohdan; Harrison, Martin James; Clase, Juha Andrew; Naylor, Neil Jason

PATENT ASSIGNEE(S): Protherics Molecular Design Limited, UK

SOURCE: PCT Int. Appl., 106 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

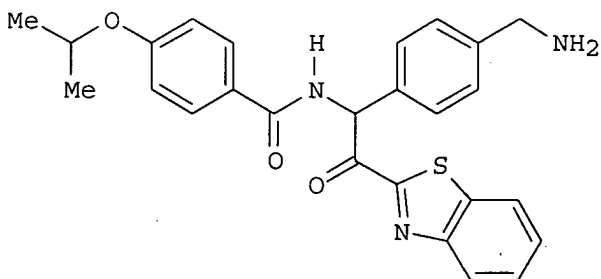
PATENT INFORMATION:

| PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|---|------|----------|-----------------|----------|
| WO 2001027096 | A1 | 20010419 | WO 2000-GB3832 | 20001005 |
| W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM | | | | |
| RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG | | | | |

PRIORITY APPLN. INFO.: GB 1999-23710 A 19991008

OTHER SOURCE(S): MARPAT 134:311433

GI



I

AB R1CHR2(CH2)aZCH2NH2 [R1 = H, NH2, NHZ1(CH2)bR3; R2 = H when R1 = NHZ1(CH2)bR3 or COR4; R3 = alk(en)yl, heterocyclyl, aryl, etc.; R4 = COR5, CF2R6, 2-(benz)oxazolyl, 2-(benz)imidazolyl, etc.; R5 = (fluoro)alkyl, alkoxy, aryl, etc.; R6 = F, (fluoro)alkyl, aryl, etc.; Z = 1,4-phenylene, 5-membered heteroarylene, etc.; Z1 = bond, CO CO2, CONH, SO2; a = 0-2; b = 0-4] were prepared as tryptase inhibitors (no data). Thus, 4-BrC6H4CH2CO2H was converted in 7 steps to 4-(BocHNH2C)C6H4CH(NH2)CO2Me which was amidated by 4-(Me2HCO)C6H4CO2H and the product condensed with

benzothiazole to give, after deprotection, title compound I.

IT 334989-05-8P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of (hetero)arylmethylamines as tryptase inhibitors)

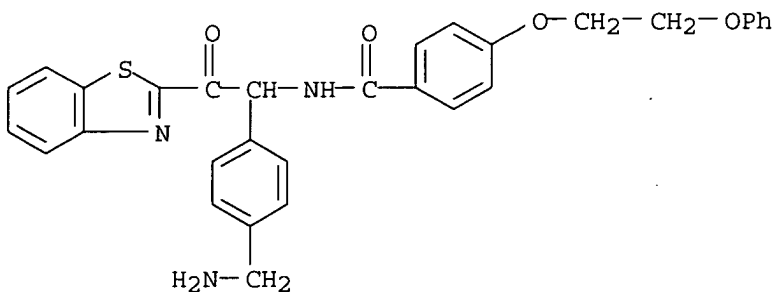
RN 334989-05-8 CAPLUS

CN Benzamide, N-[1-[4-(aminomethyl)phenyl]-2-(2-benzothiazolyl)-2-oxoethyl]-4-(2-phenoxyethoxy)-, mono(trifluoroacetate) (9CI) (CA INDEX NAME)

CM 1

CRN 334989-04-7

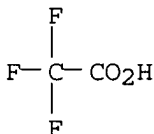
CMF C31 H27 N3 O4 S



CM 2

CRN 76-05-1

CMF C2 H F3 O2



REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L9 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2001:12407 CAPLUS

DOCUMENT NUMBER: 134:71392

TITLE: Preparation and effect of Substituted phenoxyacetic acids in complications arising from diabetes mellitus

INVENTOR(S): Van Zandt, Michael C.

PATENT ASSIGNEE(S): The Institutes for Pharmaceutical Discovery, LLC, USA

SOURCE: PCT Int. Appl., 144 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

| PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|---------------|------|----------|-----------------|----------|
| WO 2001000566 | A2 | 20010104 | WO 2000-US17377 | 20000623 |
| WO 2001000566 | A3 | 20020207 | | |

W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,
DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

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|---------------|----|----------|-----------------|----------|
| CA 2385798 | AA | 20010104 | CA 2000-2385798 | 20000623 |
| BR 2000011928 | A | 20020409 | BR 2000-11928 | 20000623 |
| EP 1198451 | A2 | 20020424 | EP 2000-944834 | 20000623 |

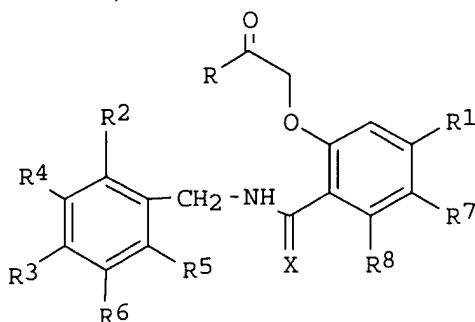
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IE, SI, LT, LV, FI, RO, MK, CY, AL

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|---------------|----|----------|----------------|----------|
| US 6420426 | B1 | 20020716 | US 2000-603817 | 20000623 |
| JP 2003503381 | T2 | 20030128 | JP 2001-506979 | 20000623 |
| EE 200100708 | A | 20030217 | EE 2001-708 | 20000623 |
| NZ 516290 | A | 20040326 | NZ 2000-516290 | 20000623 |
| NO 2001006272 | A | 20020117 | NO 2001-6272 | 20011220 |
| ZA 2002000300 | A | 20030613 | ZA 2002-300 | 20020114 |
| BG 106351 | A | 20020930 | BG 2002-106351 | 20020125 |
| US 2003036558 | A1 | 20030220 | US 2002-195964 | 20020716 |

PRIORITY APPLN. INFO.:

| | | |
|-----------------|----|----------|
| US 1999-141068P | P | 19990625 |
| US 2000-603817 | A3 | 20000623 |
| WO 2000-US17377 | W | 20000623 |

OTHER SOURCE(S): MARPAT 134:71392
GI



I

AB Disclosed are substituted phenoxyacetic acids [I; R = OH, OCH₂CH₃, C₆H₅CH₂O, (CH₃)₂CHCH₂CH₂O, CH₃(CH₂)₇O, CH₃(CH₂)₃; R₁ = H, Cl, F, Br, CH₃O, CH₃, NO₂, CH₃S, CF₃O, CH₃SO₂; R₂ = F, H; R₃ = Br, H, Cl, CH₃O, CF₃O, CH₃; R₄ = H, F, CF₃, NO₂, CH₃O, Cl; R₅ = H, F; R₆ = H, F, CF₃, CH₃O; R₇ = H, CH₃, CF₃O, NH₂, F; R₈ = H, F; X = O, S; etc] useful in the treatment of chronic complications arising from diabetes mellitus. Also disclosed are pharmaceutical compns. containing title compds. I, alone or in combination with other therapeutic agents, and methods of treatment employing the compds. and pharmaceutical compns., as well as methods for their synthesis. The pharmaceutical composition contains angiotensin converting enzyme inhibitor (benazepril, captopril, delapril, etc). Thus, title compound I (R₁ = Cl; R = OH; X = O; R₄ = NO₂; R₂, R₃, R₅, R₆, R₇, R₈ each = H) was prepared and tested.

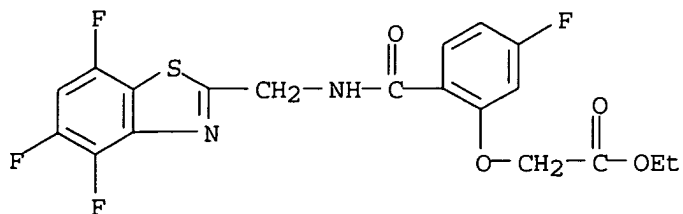
IT 314298-38-9P 314298-39-0P

RL: RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation and effect of substituted phenoxyacetic acids in complications arising from diabetes mellitus)

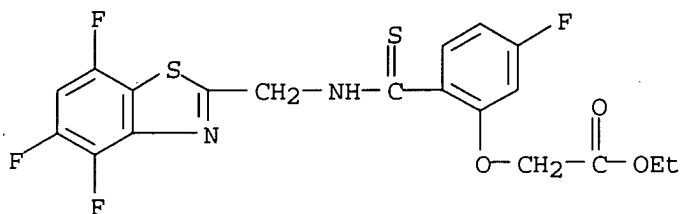
RN 314298-38-9 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]-, ethyl ester (9CI) (CA INDEX NAME)



RN 314298-39-0 CAPLUS

CN Acetic acid, [5-fluoro-2-[[thioxo[[4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]-, ethyl ester (9CI) (CA INDEX NAME)

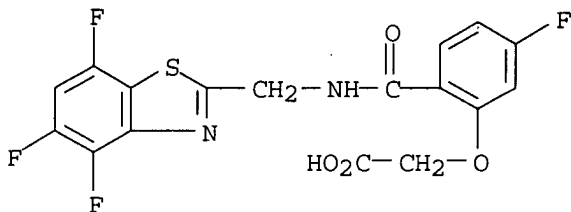


IT 314297-79-5P 314297-80-8P 314297-81-9P
314297-82-0P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation and effect of substituted phenoxyacetic acids in complications arising from diabetes mellitus)

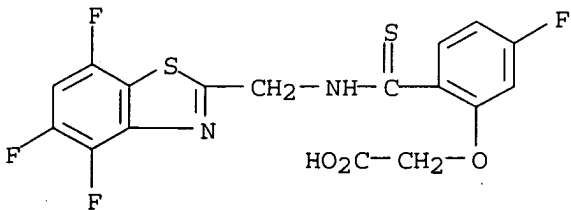
RN 314297-79-5 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



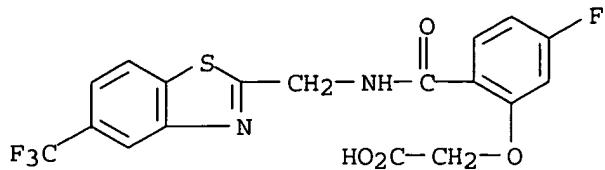
RN 314297-80-8 CAPLUS

CN Acetic acid, [5-fluoro-2-[[thioxo[[4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]- (9CI) (CA INDEX NAME)

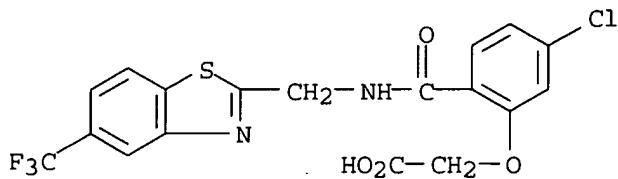


RN 314297-81-9 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[5-(trifluoromethyl)-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



RN 314297-82-0 CAPLUS
 CN Acetic acid, [5-chloro-2-[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy] - (9CI) (CA INDEX NAME)



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|--|------------|---------|
| => log y | | |
| COST IN U.S. DOLLARS | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| FULL ESTIMATED COST | 16.17 | 346.61 |
| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| CA SUBSCRIBER PRICE | -2.19 | -2.19 |

STN INTERNATIONAL LOGOFF AT 08:22:38 ON 02 MAR 2005